FILE 'REGISTRY' ENTERED AT 13:26:47 ON 27 MAY 2008 L1 STRUCTURE UPLOADED 11 S L1 157 S L1 SSS FULL L2 L3

FILE 'CAPLUS' ENTERED AT 13:27:32 ON 27 MAY 2008 L42 S L3

=> file registry COST IN U.S. DOLLARS

SINCE FILE TOTAL. ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 26 MAY 2008 HIGHEST RN 1022798-85-1 DICTIONARY FILE UPDATES: 26 MAY 2008 HIGHEST RN 1022798-85-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

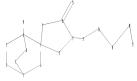
TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

Uploading C:\Program Files\Stnexp\Queries\10563271verify.str



```
13 14 15 17 18 19 20
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12
chain bonds :
```

4-14 10-13 11-15 15-18 17-20 18-19 19-20 ring bonds :

1-2 1-6 1-7 2-3 3-4 4-5 4-8 5-6 5-9 5-12 7-8 9-10 10-11 11-12 exact/norm bonds : 

exact bonds :

4-14

chain nodes :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:CLASS 14:CLASS 15:Atom 17:Atom 18:CLASS 19:CLASS

20:CLASS

Generic attributes :

15:

Saturation : Unsaturated Number of Carbon Atoms : less than 7

Number of Carbon Atoms : less than Type of Ring System : Monocyclic 17:

17:
Saturation : Unsaturated
Number of Carbon Atoms : less than 7
Type of Ring System : Monocyclic

## L1 STRUCTURE UPLOADED

=> s 11

SAMPLE SEARCH INITIATED 13:27:09 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 37 TO ITERATE

100.0% PROCESSED 37 ITERATIONS 11 ANSWERS SEARCH TIME: 00.00.01

L2 11 SEA SSS SAM L1

=> d 12 scan

L2 11 ANSMERS REGISTRY COPYRIGHT 2008 ACS on STN IN Spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one, 3'-[5-(2-thiazolvl)-3-thienyl]-, (3R)-

MF C16 H17 N3 O2 S2

Absolute stereochemistry.

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 11 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

TN Spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one, 3'-(4-pheny1-2-oxazoly1)-, (3R)-

C18 H19 N3 O3 ME

Absolute stereochemistry.

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

11 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

TN Spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one, 3'-[5-(3-pyridiny1)-2-thiazoly1]-

C17 H18 N4 02 S

MF

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s 11 sss full

FULL SEARCH INITIATED 13:27:26 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED -692 TO ITERATE

100.0% PROCESSED 692 ITERATIONS

SEARCH TIME: 00.00.01

157 ANSWERS

L3 157 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 178.36 178.57

FILE 'CAPLUS' ENTERED AT 13:27:32 ON 27 MAY 2008

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FILE COVERS 1907 - 27 May 2008 VOL 148 ISS 22 FILE LAST UPDATED: 26 May 2008 (20080526/ED)

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## => a 13

## A IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system. For a list of commands available to you in the current file, enter "HBLP COMMANDS" at an arrow prompt (=>).

=> d 14 ti abs bib

- L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN
- ${\tt TI}$   $\,$  Preparation of spiro-oxazolidinone compounds as nicotinic acetylcholine receptor ligands

AB Title compds. I [Ar1, Ar2 = 5- or 6-membered aromatic or heteroarom. moiety having 0,1 or 2 nitrogen atoms, 0 or 1 oxygen atoms, and 0 or 1 sulfur atoms; wherein Arl is unsubstituted or has 1, 2 or 3 substituents selected from alkyl, alkenyl, alkynyl, etc. and Ar2 is unsubstituted or has 1, 2 or 3 substituents selected from -CONR1R2, -NR1COR2; R1, R2 = H, alkyl, or -NR1R2 in combination is -(CH2) jG(CH2)k-; G = bond, oxygen, sulfur, etc.; j = 2-4; k = 0-2] or stereoisomers, enantiomers, in vivo hydrolysable precursors and pharmaceutically acceptable salts thereof were prepared For example, Pd(PPh3)4 catalyzed coupling reaction of 4-(N,Ndimethylaminocarbonyl)phenylboronic acid with 2.5-dibromothiophene followed by reaction with (3S)-spiro[1-azabicyclo[2.2.2]octane-3,5'oxazolidin]-2'-one afforded compound II. Compds. I are claimed useful as nicotinic acetylcholine receptor ligands for the treatment of anxiety, schizophrenia, etc. (no data).

ΆN 2006:608651 CAPLUS <<LOGINID::20080527>>

DN 145:83311

ΤI Preparation of spiro-oxazolidinone compounds as nicotinic acetylcholine receptor ligands IN Chapdelaine, Marc; Chang, Hui-Fang; Herzog, Keith J.; Horchler, Carey;

Phillips, Eifion

PA Astrazeneca AB, Swed. PCT Int. Appl., 44 pp.

SO CODEN: PIXXD2

DT Patent

English LA

FAN CNT 1

FAN.						KIN		DATE			APPLICATION NO.						DATE			
PI	WO	WO 2006065209						20060622			WO 2	2005-	SE19	E1909		2	213			
		W: AE, AG, AL,		AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,				
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			IS,	IT,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	ВJ,		
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	AU 2005317280 CA 2591430 EP 1831231																			
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		R:										ES,								
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	MX 200706743									MX 2007-6743										
	US 20080113983								IN 2007-DN4472											
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	KR 2007090922										KR 2007-713375									
	NO 2007003551															20070709				
			101124232 A 2004-636362P P							CN 2005-80048394					2	0070	815			
PRAI								2004												
		2005																		
		2005				W		2005	1213											
OS	MAI	RPAT	145:	8331	1															

RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

- L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN
- TI A preparation of derivatives of oxazolidinone with affinity to the a7-nicotinic acetylcholine receptor

- AB The invention relates to a preparation of derivs. of oxazolidinone of formula Q-X-A-Y (wherein: Q is spiro(azabicyclooctanoxazolidinone) derivative; A is O, S, or NH, etc.; X is 5- or 6-membered heterocycle; Y is is 5- or 6-membered (hetero) aromatic ringl with affinity to the  $\alpha 7-\text{mictorinic}$  acetylcholine receptor. For instance, oxazolidinone derivative I was prepared via phenylation of II by phenylboronic acid. The compds. of the invention were screened in  $\alpha 7$  nAChR subtype affinity assay and showed binding affinities (Ki) of less than 1000 nM.
- AN 2005:58211 CAPLUS <<LOGINID::20080527>>
- DN 142:155977
- TI A preparation of derivatives of oxazolidinone with affinity to the  $\alpha^7$ -nicotinic acetylcholine receptor
- IN Chang, Hui-Fang; Phillips, Eifion
- PA Astrazeneca AB, Swed.; Astrazeneca UK Limited
- SO PCT Int. Appl., 77 pp. CODEN: PIXXD2
- DT Patent
- LA English
- FAN.CNT 1

PATENT NO. KIND	T, AU, AZ,	APPLICATION NO. WO 2004-GB2904 BA, BB, BG, BR, BW, BY	20040706		
PI WO 2005005435 A1	T, AU, AZ,				
	T, AU, AZ,				
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LK, LR, LS, LT, I	U, LV, MA,	MD, MG, MK, MN, MW, MX	MZ, NA, NI,		
NO, NZ, OM, PG, F	H, PL, PT,	RO, RU, SC, SD, SE, SG	, SK, SL, SY,		
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AU 2004255920 A1	20050120	20040706			

	CA	25315	10			A1		2005	0120	(	CA	2004-	2531	510		2	0040	706
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	NO	20060	006	12		A		2006	0406	ľ	10	2006-	612			2	0060	208
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	WO	2004-	GB29	904		W		2004	0706									
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